L Number	Hits	Search Text	DB	Time stamp
1	0	("fluoren$").PN.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
2	28987	fluoren$	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
3	2024	RAR	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
4	5257	fluorenone	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
5	0	RAR and fluorenone	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
6	0	9807716.pn	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
7	586	562/466.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
8	189	549/26.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
9	108	514/437.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
10	524	514/569.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
11	1331	562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
12	6541	retinoid	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
13	26450	phenanth$	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
14	8441	retinoic	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
15	113	retinoic and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
16	853	560/56.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
17	331	514/544.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42

18	2	5945561.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
19	2	6319948.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
20	2	5760084.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
21	2	5624957.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
22	45	fluoren$ and RAR	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
23	5	9807716.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42
24	99	arotinoid	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
25	1	arotinoid and 562/466.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
26	2	5618839.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
27	17	5618839.URPN.	USPAT	2004/04/21 05:43
28	3	arotinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
29	139	retinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
30	4	(retinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)) and phenanth$	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
31	3	phenanth$ and (retinoic and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.))	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
32	2	5075487.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43
33	853	560/56.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 06:35
34	331	514/544.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 06:35

	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
1	IS&R	L1	0	("fluoren$").PN.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
2	BRS	L2	28987	fluoren$	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
3	BRS	L3	2024	RAR	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
4	BRS	L4	5257	fluorenone	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
5	BRS	L5	0	RAR and fluorenone	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
6	BRS	L6	0	9807716.pn	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
7	BRS	L7	586	562/466.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		

	Errors
1	0
2	0
3	0
4	0
5	0
6	0
7	0

	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
8	BRS	L8	189	549/26.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
9	BRS	L9	108	514/437.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
10	BRS	L10	524	514/569.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
11	BRS	L11	1331	562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
12	BRS	L12	6541	retinoid	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
13	BRS	L13	26450	phenanth$	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
14	BRS	L14	8441	retinoic	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		

	Errors
8	0
9	0
10	0
11	0
12	0
13	0
14	0

	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
15	BRS	L15	113	retinoic and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
16	BRS	L16	853	560/56.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
17	BRS	L17	331	514/544.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
18	BRS	L18	2	5945561.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
19	BRS	L19	2	6319948.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
20	BRS	L20	2	5760084.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
21	BRS	L21	2	5624957.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		

	Errors
15	0
16	0
17	0
18	0
19	0
20	0
21	0

	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
22	BRS	L22	45	fluoren$ and RAR	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
23	BRS	L23	5	9807716.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:42		
24	BRS	L24	99	arotinoid	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
25	BRS	L25	1	arotinoid and 562/466.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
26	BRS	L26	2	5618839.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
27	BRS	L27	17	5618839.URPN.	USPAT	2004/04/21 05:43		
28	BRS	L28	3	arotinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
29	BRS	L29	139	retinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		

	Errors
22	0
23	0
24	0
25	0
26	0
27	0
28	0
29	0

	Type	L #	Hits	Search Text	DBs	Time Stamp	Comments	Error Definition
30	BRS	L30	4	(retinoid and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.)) and phenanth$	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
31	BRS	L31	3	phenanth$ and (retinoic and (562/466.ccls. or 549/26.ccls. or 514/437.ccls. or 514/569.ccls.))	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
32	BRS	L32	2	5075487.pn.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 05:43		
33	BRS	L33	853	560/56.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 06:35		
34	BRS	L34	331	514/544.ccls.	USPAT; US-PGPUB; EPO; JPO; DERWENT	2004/04/21 06:35		

	Errors
30	0
31	0
32	0
33	0
34	0

```
Welcome to STN International!  Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:
TERMINAL (ENTER 1, 2, 3, OR ?):2


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 NEWS  1          Web Page URLs for STN Seminar Schedule - N. America
 NEWS  2          "Ask CAS" for self-help around the clock
 NEWS  3  JAN 27  Source of Registration (SR) information in REGISTRY updated
                  and searchable
 NEWS  4  JAN 27  A new search aid, the Company Name Thesaurus, available in
                  CA/CAplus
 NEWS  5  FEB 05  German (DE) application and patent publication number format
                  changes
 NEWS  6  MAR 03  MEDLINE and LMEDLINE reloaded
 NEWS  7  MAR 03  MEDLINE file segment of TOXCENTER reloaded
 NEWS  8  MAR 03  FRANCEPAT now available on STN
 NEWS  9  MAR 29  Pharmaceutical Substances (PS) now available on STN
 NEWS 10  MAR 29  WPIFV now available on STN
 NEWS 11  MAR 29  No connect hour charges in WPIFV until May 1, 2004
 NEWS 12  MAR 29  New monthly current-awareness alert (SDI) frequency in RAPRA

 NEWS EXPRESS  MARCH 31 CURRENT WINDOWS VERSION IS V7.00A, CURRENT
               MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),
               AND CURRENT DISCOVER FILE IS DATED 13 APRIL 2004
 NEWS HOURS    STN Operating Hours Plus Help Desk Availability
 NEWS INTER    General Internet Information
 NEWS LOGIN    Welcome Banner and News Items
 NEWS PHONE    Direct Dial and Telecommunication Network Access to STN
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Enter NEWS followed by the item number or name to see news on that
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FILE 'HOME' ENTERED AT 05:54:00 ON 21 APR 2004

=> FIL STNGUIDE
COST IN U.S. DOLLARS                                SINCE FILE      TOTAL
                                                      ENTRY    SESSION
FULL ESTIMATED COST                                    0.21       0.21

FILE 'STNGUIDE' ENTERED AT 05:54:06 ON 21 APR 2004
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AND TECHNOLOGY CORPORATION, AND FACHINFORMATIONSZENTRUM KARLSRUHE

FILE CONTAINS CURRENT INFORMATION.
LAST RELOADED: Apr 16, 2004 (20040416/UP).
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=> FIL HOME
COST IN U.S. DOLLARS                                    SINCE FILE      TOTAL
                                                           ENTRY    SESSION
FULL ESTIMATED COST                                         0.06       0.27

FILE 'HOME' ENTERED AT 05:54:10 ON 21 APR 2004

=> file reg
COST IN U.S. DOLLARS                                    SINCE FILE      TOTAL
                                                           ENTRY    SESSION
FULL ESTIMATED COST                                         0.21       0.48

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STRUCTURE FILE UPDATES:   19 APR 2004  HIGHEST RN 676225-08-4
DICTIONARY FILE UPDATES:  19 APR 2004  HIGHEST RN 676225-08-4

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

  Please note that search-term pricing does apply when
  conducting SmartSELECT searches.

Crossover limits have been increased.  See HELP CROSSOVER for details.

Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html

=>
Uploading C:\Examination Auxillary files\10075845\10075845 generic six allowance.str
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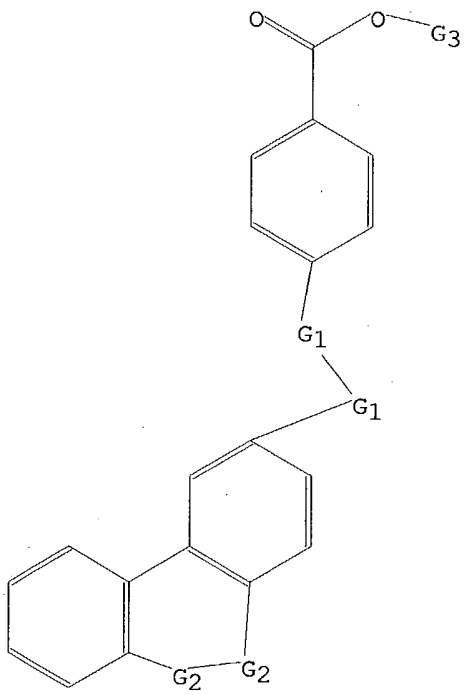
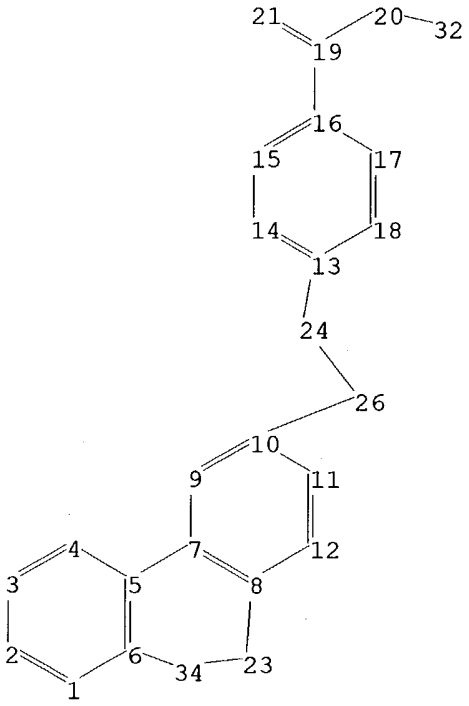
Ak*1



28*1



chain nodes :
19 20 21 24 26 28 32
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 23 34
chain bonds :
10-26 13-24 16-19 19-20 19-21 20-32 24-26
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-34 7-8 7-9 8-12 8-23 9-10 10-11 11-12
13-18 13-14 14-15 15-16 16-17 17-18 23-34
exact/norm bonds :
5-7 6-34 8-23 10-26 13-24 16-19 19-20 19-21 20-32 23-34 24-26
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-9 8-12 9-10 10-11 11-12 13-18 13-14
14-15
15-16 16-17 17-18

G1:C,O,S,N

G2:C,O,S

G3:H,[*1]

Hydrogen count :
1:>= minimum 1 2:>= minimum 1 3:>= minimum 1 4:>= minimum 1
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:CLASS
20:CLASS 21:CLASS
23:CLASS 24:CLASS 26:CLASS 28:CLASS 32:CLASS 34:CLASS
Element Count :

```
Node 28: Limited
    C,C1-7



L1      STRUCTURE UPLOADED

=> d l1
L1 HAS NO ANSWERS
L1              STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY -  AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.


=> search l1 sss sam
SAMPLE SEARCH INITIATED 05:54:49 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED -   3439 TO ITERATE

 29.1% PROCESSED    1000 ITERATIONS                                  0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS:  ONLINE  **COMPLETE**
                        BATCH   **COMPLETE**
PROJECTED ITERATIONS:        65264 TO    72296
PROJECTED ANSWERS:               0 TO        0

L2            0 SEA SSS SAM L1

=> search l1 sss full
FULL SEARCH INITIATED 05:55:01 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED -  69812 TO ITERATE

100.0% PROCESSED   69812 ITERATIONS                               0 ANSWERS
SEARCH TIME: 00.00.03

L3            0 SEA SSS FUL L1

=>
Uploading C:\Examination Auxillary files\10075845\10075845 generic fiveallowance.str
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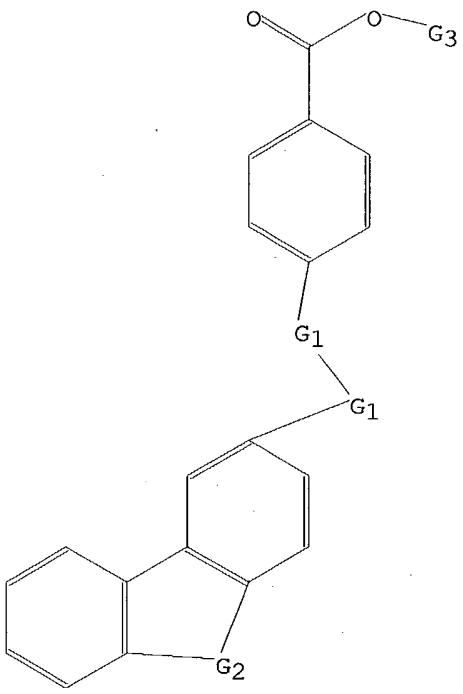
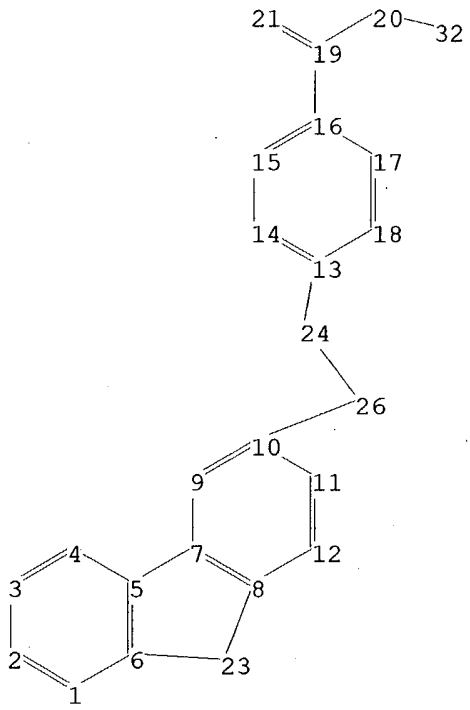
Ak*1



28*1



chain nodes :
19 20 21 24 26 28 32
ring nodes :
1 2 3 4 5 6 7 8 9 10 11 12 13 14 15 16 17 18 23
chain bonds :
10-26 13-24 16-19 19-20 19-21 20-32 24-26
ring bonds :
1-2 1-6 2-3 3-4 4-5 5-6 5-7 6-23 7-8 7-9 8-12 8-23 9-10 10-11 11-12
13-18 13-14 14-15 15-16 16-17 17-18
exact/norm bonds :
5-7 6-23 8-23 10-26 13-24 16-19 19-20 19-21 20-32 24-26
normalized bonds :
1-2 1-6 2-3 3-4 4-5 5-6 7-8 7-9 8-12 9-10 10-11 11-12 13-18 13-14
14-15
15-16 16-17 17-18

G1:C,O,S,N

G2:C,O,S

G3:H,[*1]

Hydrogen count :
1:>= minimum 1 2:>= minimum 1 3:>= minimum 1 4:>= minimum 1
Match level :
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:Atom 8:Atom 9:Atom 10:Atom
11:Atom 12:Atom 13:Atom 14:Atom 15:Atom 16:Atom 17:Atom 18:Atom 19:CLASS
20:CLASS 21:CLASS
23:CLASS 24:CLASS 26:CLASS 28:CLASS 32:CLASS
Element Count :

Node 28: Limited
C,C1-7

L4 STRUCTURE UPLOADED

=> d l4
L4 HAS NO ANSWERS
L4 STR

* STRUCTURE DIAGRAM TOO LARGE FOR DISPLAY - AVAILABLE VIA OFFLINE PRINT *

Structure attributes must be viewed using STN Express query preparation.

=> search l4 sss sam
SAMPLE SEARCH INITIATED 05:57:01 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 3288 TO ITERATE

30.4% PROCESSED 1000 ITERATIONS 0 ANSWERS
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
BATCH **COMPLETE**
PROJECTED ITERATIONS: 62322 TO 69198
PROJECTED ANSWERS: 0 TO 0

L5 0 SEA SSS SAM L4

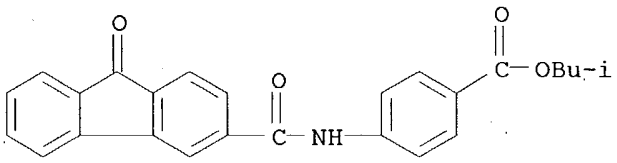
=> search l4 sss full
FULL SEARCH INITIATED 05:57:09 FILE 'REGISTRY'
FULL SCREEN SEARCH COMPLETED - 66736 TO ITERATE

100.0% PROCESSED 66736 ITERATIONS 3 ANSWERS
SEARCH TIME: 00.00.03

L6 3 SEA SSS FUL L4

=> d scan

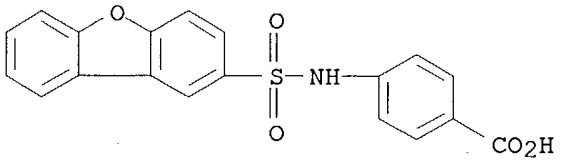
L6 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Benzoic acid, 4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl ester (9CI)
MF C25 H21 N O4



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

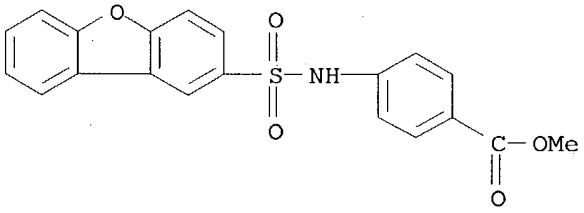
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):3

L6 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Benzoic acid, 4-[(2-dibenzofuranylsulfonyl)amino]- (9CI)
MF C19 H13 N O5 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

L6 3 ANSWERS REGISTRY COPYRIGHT 2004 ACS on STN
IN Benzoic acid, 4-[(2-dibenzofuranylsulfonyl)amino]-, methyl ester (9CI)
MF C20 H15 N O5 S



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

ALL ANSWERS HAVE BEEN SCANNED

=> file caplus
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	312.94	313.42

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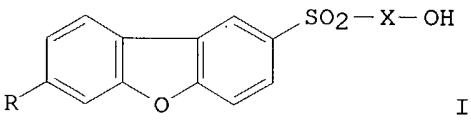
FILE LAST UPDATED: 20 Apr 2004 (20040420/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> l6
L7 1 L6

=> d l7 ti fbib abs

L7 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2004 ACS on STN
TI Synthesis and biological activity of some new dibenzofuran- and 7-nitrodibenzofuran-2-sulfonyl amino acid derivatives
AN 1985:505304 CAPLUS
DN 103:105304
TI Synthesis and biological activity of some new dibenzofuran- and 7-nitrodibenzofuran-2-sulfonyl amino acid derivatives
AU El-Naggar, A. M.; Abd El-Salam, A. M.; Ahmed, F. S. M.; Ibrahim, T. M.
CS Fac. Sci., Al-Azhar Univ., Nasr, Egypt
SO Acta Pharmaceutica Jugoslavica (1985), 35(1), 15-22
CODEN: APJUA8; ISSN: 0001-6667
DT Journal
LA English
GI



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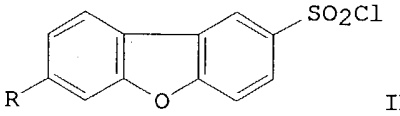

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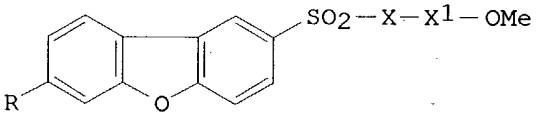

III

AB Title amino acid derivs. I (X = β-Ala, Val, Leu, p-NHC6H4CO, Phe, etc.; R = H or NO2) were prepared by sulfonylating the corresponding amino acid with sulfonyl chlorides II (R = H or NO2). I were esterified with MeOH via SOCl2 to give the corresponding Me esters. Also, I were coupled with amino acid Me ester hydrochlorides by DCC in THF containing Et3N to give the corresponding dipeptides, e.g. III (X-X1 = DL-Val-DL-Val, Pro-Phe, R = H; X-X1 = Pro-DL-Ser, Leu-Tyr, R = NO2). Nineteen synthesized compds., e.g. I (X = Leu, R = H; X = β-Ala, R = NO2) and III (X-X1 = Tyr-Phe, R = NO2), were active against various microorganisms, e.g. Bacillus subtilis or B. cereus.

=> logoff hold
COST IN U.S. DOLLARS

	SINCE FILE ENTRY	TOTAL SESSION
FULL ESTIMATED COST	3.86	317.28

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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)        SINCE FILE      TOTAL
                                                     ENTRY    SESSION
CA SUBSCRIBER PRICE                                  -0.69      -0.69

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PASSWORD:
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FULL ESTIMATED COST                                   3.86     317.28

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CA SUBSCRIBER PRICE                                  -0.69      -0.69

=> file reg
COST IN U.S. DOLLARS                              SINCE FILE      TOTAL
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FULL ESTIMATED COST                                   4.29     317.71

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DICTIONARY FILE UPDATES:  19 APR 2004  HIGHEST RN 676225-08-4

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

  Please note that search-term pricing does apply when
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Experimental and calculated property data are now available. For more
information enter HELP PROP at an arrow prompt in the file or refer
to the file summary sheet on the web at:
http://www.cas.org/ONLINE/DBSS/registryss.html
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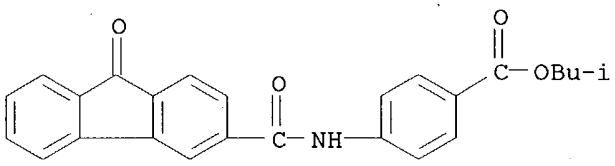
```
=> eBenzoic acid, 4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl
ester/cn
MISSING OPERATOR '4-[[(9-OXO-9H-F'

=> e Benzoic acid, 4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl
ester/cn
E1            1      BENZOIC ACID, 4,6-DIMETHOXY-3-(5-METHOXY-2-(METHOXYCARBONYL)
                     -3-PENTYLPHENOXY)-2-PENTYL-, METHYL ESTER/CN
E2            1      BENZOIC ACID, 4,6-DIMETHOXY-3-METHYL-2-(1-OXOBUTOXY)-, METHY
                     L ESTER/CN
E3            0 --> BENZOIC ACID, 4-(9-OXO-9H-FLUOREN-3-YL)CARBONYL AMINO-, 2
                     -METHYLPROPYL     ESTER/CN
E4            1      BENZOIC ACID, 4-(((((((2,3-DIHYDRO-2,2-DIMETHYL-7-BENZOFURA
                     NYL)OXY)CARBONYL)METHYLAMINO)CARBONYL)AMINO)SULFONYL)AMINO)-
                     , ETHYL ESTER/CN
E5            1      BENZOIC ACID, 4-((((((2,3,4-TRICHLOROPHENYL)THIO)ACETYL)OXY
                     )ACETYL)AMINO)METHYL)-, METHYL ESTER/CN
E6            1      BENZOIC ACID, 4-((((((3B)-17-OXOANDROST-5-EN-3-YL)AMIN
                     O)CARBONYL)OXY)METHYL)AMINO)-, 2-(DIETHYLAMINO)ETHYL ESTER/C
                     N
E7            1      BENZOIC ACID, 4-((((((3B)-17-OXOANDROST-5-EN-3-YL)OXY)
                     METHYL)THIO)METHYL)AMINO)-, 2-(DIMETHYLAMINO)ETHYL ESTER/CN
E8            1      BENZOIC ACID, 4-((((((3B,17B)-3-HYDROXYANDROST-5-
                     EN-17-YL)AMINO)CARBONYL)OXY)METHYL)AMINO)-, 2-(DIETHYLAMINO)
                     ETHYL ESTER/CN
E9            1      BENZOIC ACID, 4-((((((4-BROMO-8-CHLORO-1-NAPHTHALENYL)THIO)
                     ACETYL)OXY)ACETYL)AMINO)METHYL)-, METHYL ESTER/CN
E10           1      BENZOIC ACID, 4-((((((4-CHLOROPHENYL)AMINO)CARBONYL)HYDRAZO
                     NO)PHENYLMETHYL)THIO)METHYL)-, 1,1-DIMETHYLETHYL ESTER/CN
E11           1      BENZOIC ACID, 4-((((((4-CHLOROPHENYL)SULFONYL)AMINO)PHENYLM
                     ETHYLENE)AMINO)THIOXOMETHYL)AMINO)-/CN
E12           1      BENZOIC ACID, 4-((((((4-METHOXYPHENYL)SULFONYL)(PHENYLMETHY
                     L)AMINO)ACETYL)AMINO)OXY)DIPHENYLMETHYL)-/CN

=> e Benzoic acid, 4-(((9-oxo-9H-fluoren-3-yl)carbonyl)amino)-,2-methylpropyl
ester/cn
E1            1      BENZOIC ACID, 4-(((9-METHYL-4-OXO-2-(1-PYRROLIDINYL)-4H-PYRI
                     DO(1,2-A)PYRIMIDIN-3-YL)METHYLENE)HYDRAZINO)-/CN
E2            1      BENZOIC ACID, 4-(((9-OXO-9H-FLUOREN-3-YL)CARBONYL)AMINO)-, 2
                     -METHYLPROPYL ESTER/CN
E3            0 --> BENZOIC ACID, 4-(((9-OXO-9H-FLUOREN-3-YL)CARBONYL)AMINO)-,2-
                     METHYLPROPYL ESTER/CN
E4            1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-/C
                     N
E5            1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-,
                     1-(4-(2-AMINO-2-OXOETHOXY)-5-METHOXY-2-NITROPHENYL)ETHYL EST
                     ER/CN
E6            1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-,
                     1-METHYLETHYL ESTER/CN
E7            1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-2-
                     HYDROXY-/CN
E8            1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-2-
                     HYDROXY-, METHYL ESTER/CN
E9            1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-2-
                     NITRO-/CN
E10           1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-3-
                     IODO-/CN
E11           1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-3-
                     IODO-, METHYL ESTER/CN
E12           1      BENZOIC ACID, 4-(((9H-FLUOREN-9-YLMETHOXY)CARBONYL)AMINO)-3-
                     METHOXY-/CN
```

=> e2
L8 1 "BENZOIC ACID, 4-(((9-OXO-9H-FLUOREN-3-YL)CARBONYL)AMINO)-, 2-METHYLPROPYL ESTER"/CN

=> d l8

L8 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2004 ACS on STN
RN 505063-62-7 REGISTRY
CN **Benzoic acid, 4-[[(9-oxo-9H-fluoren-3-yl)carbonyl]amino]-, 2-methylpropyl ester (9CI)** (CA INDEX NAME)
FS 3D CONCORD
MF C25 H21 N O4
SR Chemical Library
LC STN Files: CHEMCATS



PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

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DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
CA SUBSCRIBER PRICE	0.00	-0.69

SESSION WILL BE HELD FOR 60 MINUTES
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